

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021622\
 Data File : BP009175.D
 Acq On : 16 Feb 2022 13:01
 Operator : CG/JU
 Sample : PB142605BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB142605BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 02/17/2022
 Supervised By : Jagrut Upadhyay 02/17/2022

Quant Time: Feb 16 14:32:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 16 14:30:24 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	227289	20.000	ng	0.00
21) Naphthalene-d8	10.581	136	943505	20.000	ng	0.00
39) Acenaphthene-d10	14.428	164	672808	20.000	ng	0.00
64) Phenanthrene-d10	17.186	188	1457071	20.000	ng	0.00
76) Chrysene-d12	21.286	240	1502395	20.000	ng	0.00
86) Perylene-d12	23.674	264	1546414	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	1660009	130.281	ng	0.00
7) Phenol-d6	6.981	99	2171862	129.359	ng	0.00
23) Nitrobenzene-d5	8.946	82	1294934	77.399	ng	0.00
42) 2,4,6-Tribromophenol	15.934	330	1208603	134.540	ng	0.00
45) 2-Fluorobiphenyl	13.046	172	3604301	74.994	ng	0.00
79) Terphenyl-d14	19.751	244	5988913	71.686	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.281	88	135985	32.406	ng	99
3) Pyridine	3.681	79	353391	28.828	ng	98
4) n-Nitrosodimethylamine	3.593	42	181106	38.565	ng	# 92
6) Aniline	7.116	93	705281	37.009	ng	99
8) 2-Chlorophenol	7.352	128	673584	46.842	ng	97
9) Benzaldehyde	6.928	77	313708	36.888	ng	98
10) Phenol	7.005	94	812885	48.155	ng	96
11) bis(2-Chloroethyl)ether	7.211	93	568520	43.374	ng	98
12) 1,3-Dichlorobenzene	7.669	146	670421	40.312	ng	98
13) 1,4-Dichlorobenzene	7.816	146	697735	41.097	ng	99
14) 1,2-Dichlorobenzene	8.134	146	678344	41.229	ng	98
15) Benzyl Alcohol	8.034	79	550014	51.462	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.310	45	637539	43.154	ng	97
17) 2-Methylphenol	8.246	107	546149	48.260	ng	98
18) Hexachloroethane	8.852	117	239053	42.505	ng	96
19) n-Nitroso-di-n-propyla...	8.599	70	460910	47.946	ng	99
20) 3+4-Methylphenols	8.581	107	745329	48.122	ng	96
22) Acetophenone	8.610	105	951490	43.139	ng	# 97
24) Nitrobenzene	8.987	77	660071	41.735	ng	95
25) Isophorone	9.516	82	1253691	45.100	ng	97
26) 2-Nitrophenol	9.705	139	371986	45.475	ng	96
27) 2,4-Dimethylphenol	9.775	122	630290	51.531	ng	98
28) bis(2-Chloroethoxy)met...	9.993	93	786949	41.708	ng	99
29) 2,4-Dichlorophenol	10.257	162	718760	47.027	ng	99
30) 1,2,4-Trichlorobenzene	10.446	180	734486	40.871	ng	99
31) Naphthalene	10.628	128	1976930	41.081	ng	99
32) Benzoic acid	10.005	122	428804	43.373	ng	99
33) 4-Chloroaniline	10.763	127	493592	25.402	ng	99
34) Hexachlorobutadiene	10.910	225	457470	41.382	ng	99
35) Caprolactam	11.575	113	219117m	48.892	ng	
36) 4-Chloro-3-methylphenol	11.904	107	691382	47.421	ng	99
37) 2-Methylnaphthalene	12.246	142	1531941	44.986	ng	97
38) 1-Methylnaphthalene	12.463	142	1417595	42.602	ng	97
40) 1,2,4,5-Tetrachloroben...	12.622	216	913938	42.676	ng	99
41) Hexachlorocyclopentadiene	12.593	237	535401	68.262	ng	98
43) 2,4,6-Trichlorophenol	12.869	196	616527	43.413	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021622\
 Data File : BP009175.D
 Acq On : 16 Feb 2022 13:01
 Operator : CG/JU
 Sample : PB142605BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB142605BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 02/17/2022
 Supervised By : Jagrut Upadhyay 02/17/2022

Quant Time: Feb 16 14:32:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 16 14:30:24 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.963	196	671930	44.121	ng	98
46) 1,1'-Biphenyl	13.257	154	2066083	41.796	ng	99
47) 2-Chloronaphthalene	13.304	162	1572788	40.040	ng	99
48) 2-Nitroaniline	13.522	65	406907	44.707	ng	98
49) Acenaphthylene	14.151	152	2503230	42.270	ng	100
50) Dimethylphthalate	13.893	163	2074559	43.257	ng	100
51) 2,6-Dinitrotoluene	14.016	165	465118	46.501	ng	93
52) Acenaphthene	14.493	154	1473615	40.897	ng	98
53) 3-Nitroaniline	14.351	138	311476	30.445	ng	97
54) 2,4-Dinitrophenol	14.581	184	512744	76.610	ng	95
55) Dibenzofuran	14.828	168	2476664	40.798	ng	98
56) 4-Nitrophenol	14.698	139	649288	75.055	ng	90
57) 2,4-Dinitrotoluene	14.822	165	655277	50.164	ng	# 82
58) Fluorene	15.481	166	1994766	42.771	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.069	232	590917m	44.338	ng	
60) Diethylphthalate	15.257	149	2043510	45.063	ng	100
61) 4-Chlorophenyl-phenyle...	15.475	204	1085252	42.828	ng	99
62) 4-Nitroaniline	15.528	138	418398	40.674	ng	91
63) Azobenzene	15.769	77	1677977	42.767	ng	95
65) 4,6-Dinitro-2-methylph...	15.598	198	352728	39.447	ng	97
66) n-Nitrosodiphenylamine	15.692	169	1808730	40.931	ng	98
67) 4-Bromophenyl-phenylether	16.369	248	713949	42.283	ng	99
68) Hexachlorobenzene	16.498	284	820677	43.390	ng	99
69) Atrazine	16.657	200	706096	48.189	ng	98
70) Pentachlorophenol	16.851	266	904338	90.807	ng	99
71) Phenanthrene	17.228	178	3246196	41.145	ng	100
72) Anthracene	17.322	178	3352972	43.438	ng	99
73) Carbazole	17.604	167	3018198	40.792	ng	99
74) Di-n-butylphthalate	18.145	149	3615790	48.231	ng	99
75) Fluoranthene	19.204	202	3989455	45.117	ng	97
77) Benzidine	19.386	184	1857210	59.158	ng	98
78) Pyrene	19.557	202	4076270	38.141	ng	99
80) Butylbenzylphthalate	20.427	149	1674901	45.577	ng	97
81) Benzo(a)anthracene	21.275	228	3921760	40.507	ng	98
82) 3,3'-Dichlorobenzidine	21.204	252	1304030	38.846	ng	98
83) Chrysene	21.322	228	3835345	41.093	ng	98
84) Bis(2-ethylhexyl)phtha...	21.186	149	2446775	50.334	ng	97
85) Di-n-octyl phthalate	22.104	149	4286431	55.124	ng	100
87) Indeno(1,2,3-cd)pyrene	26.180	276	4806850	41.834	ng	97
88) Benzo(b)fluoranthene	22.951	252	4176030	43.866	ng	98
89) Benzo(k)fluoranthene	22.998	252	4181420	44.025	ng	98
90) Benzo(a)pyrene	23.574	252	4098112	51.633	ng	# 97
91) Dibenzo(a,h)anthracene	26.186	278	4098075	43.400	ng	98
92) Benzo(g,h,i)perylene	26.939	276	4010290	42.655	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

